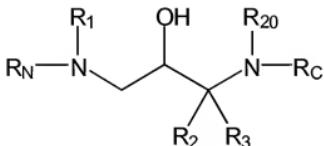


Listing of claims:

The following listing of claims will replace all prior versions and listings of claims in the application.

1. (Currently Amended) A compound of the formula I:



or a pharmaceutically acceptable salt or ester thereof, wherein R₂₀ is H, C₁₋₆ alkyl or alkenyl, C₁₋₆ haloalkyl or C₄₋₆ cycloalkyl; R₁ is -(CH₂)₂₋₄-S(O)₂₋(C₁₋₆ alkyl), or

C₁₋₆ alkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, -OH, -SH, -C≡N, -CF₃, C₁₋₆ alkoxy, amino, mono- or dialkylamino, -N(R)C(O)R', -OC(-O)-amino and -OC(-O)-mono- or dialkylamino, or

C₂₋₆ alkenyl or C₂₋₆ alkynyl, each of which is optionally substituted with 1, 2, or 3 groups independently selected from halogen, -OH, -SH, -C≡N, -CF₃, C₁₋₆ alkoxy, amino, and mono- or dialkylamino, or

aryl, heteroaryl, heterocyclyl, -C₁₋₆ alkyl-aryl, -C₁₋₆ alkyl-heteroaryl, or -C₁₋₆ alkyl-heterocyclyl, where the ring portions of each are optionally substituted with 1, 2, 3, or 4 groups independently selected from halogen, -OH, -SH, -C≡N, -NR₁₀₅R'₁₀₅, -CO₂R, -N(R)COR', or -N(R)SO₂R', -C(=O)-(C_{1-C4}) alkyl, -SO₂-amino, -SO₂-mono or dialkylamino, -C(=O)-amino, -C(=O)-mono or dialkylamino, -SO₂-(C_{1-C4}) alkyl, or

C_1-C_6 alkoxy optionally substituted with 1, 2, or 3 groups which are independently selected from halogen, or
 C_6-C_{14} cycloalkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, $-OH$, $-SH$, $-C\equiv N$, $-CF_3$, C_6-C_8 alkoxy, amino, $-C_6-C_8$ alkyl and mono- or dialkylamino, or
 C_6-C_{14} alkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, $-OH$, $-SH$, $-C\equiv N$, $-CF_3$, C_6-C_8 alkoxy, amino, mono- or dialkylamino and C_6-C_8 alkyl, or
 C_6-C_{14} alkenyl or C_6-C_{14} alkynyl each of which is optionally substituted with 1, 2, or 3 groups independently selected from halogen, $-OH$, $-SH$, $-C\equiv N$, $-CF_3$, C_6-C_8 alkoxy, amino, C_6-C_8 alkyl and mono- or dialkylamino, and the heteroacyethyl group is optionally further substituted with α -ole;

R and R' independently are hydrogen, C_1-C_{10} alkyl, C_1-C_{10} alkylaryl or C_1-C_{10} alkylheteroaryl;

R_c is hydrogen, $-(CR_{245}R_{250})_{0-4}$ -aryl, $-(CR_{245}R_{250})_{0-4}$ -heteroaryl, $-(CR_{245}R_{250})_{0-4}$ -heterocyclyl, $-(CR_{245}R_{250})_{0-4}$ -aryl-heteroaryl, $-(CR_{245}R_{250})_{0-4}$ -aryl-heterocyclyl, $-(CR_{245}R_{250})_{0-4}$ -aryl-aryl, $-(CR_{245}R_{250})_{0-4}$ -heteroaryl-aryl, $-(CR_{245}R_{250})_{0-4}$ -heteroaryl-heterocyclyl, $-(CR_{245}R_{250})_{0-4}$ -heteroaryl-heteroaryl, $-(CR_{245}R_{250})_{0-4}$ -heterocyclyl-heteroaryl, $-(CR_{245}R_{250})_{0-4}$ -heterocyclyl-aryl, $-[C(R_{255})(R_{260})]_{1-3}-CO-N-(R_{255})_2$, $-CH(aryl)_2$, $-CH(heteroaryl)_2$, $-CH(heterocyclyl)_2$, $-CH(aryl)(heteroaryl)$, $-(CH_2)_{0-1}$ - $CH((CH_2)_{0-6}-OH)-(CH_2)_{0-1}$ -aryl, $-(CH_2)_{0-1}-CH((CH_2)_{0-6}-OH)-(CH_2)_{0-1}$ -heteroaryl, $-CH(aryl$ or $-heteroaryl)-CO-O(C_1-C_4$ alkyl), $-CH(-CH_2-OH)-CH(OH)-phenyl-NO_2$, $(C_1-C_6$ alkyl) $-O-(C_1-C_6$ alkyl) $-OH$; $-CH_2-NH-CH_2-CH(-O-CH_2-CH_3)_2$, $-(CH_2)_{0-6}-C(=NR_{235})(NR_{235}R_{240})$, or

~~C₁-C₄ alkyl optionally substituted with 1, 2, or 3 groups independently selected from the group consisting of R₂₀₅, -OC=ONR₂₃₅R₂₄₀, -S(=O)₀₋₂(C₁-C₆ alkyl), -SH, -NR₂₃₅C=ONR₂₃₅R₂₄₀, -C=ONR₂₃₅R₂₄₀, and -S(=O)₀₋₂NR₂₃₅R₂₄₀, or -(CH₂)₀₋₂-(C₂-C₆) cycloalkyl wherein the cycloalkyl is optionally substituted with 1, 2, or 3 groups independently selected from the group consisting of R₂₀₅, CO₂H, and CO₂-(C₁-C₄ alkyl), or cyclopentyl, cyclohexyl, or cycloheptyl ring fused to aryl, heteroaryl, or heterocyclyl wherein one, two or three carbons of the cyclopentyl, cyclohexyl, or cycloheptyl is optionally replaced with a heteroatom independently selected from NH, NR₂₁₅, O, or S(=O)₀₋₂, and wherein the cyclopentyl, cyclohexyl, or cycloheptyl group can be optionally substituted with one or two groups that are independently R₂₀₅, =O, -CO-NR₂₃₅R₂₄₀, or -SO₂-(C₁-C₄ alkyl), or C₂-C₆ alkenyl or C₂-C₆ alkynyl, each of which is optionally substituted with 1, 2, or 3 R₂₀₅ groups, wherein each aryl and heteroaryl is optionally substituted with 1, 2, or 3 R₂₀₀, and wherein each heterocyclyl is optionally substituted with 1, 2, 3, or 4 R₂₁₀;~~

R₂₀₀ at each occurrence is independently selected from -OH, -NO₂, halogen, -CO₂H, C≡N, -(CH₂)₀₋₄-CO-NR₂₂₀R₂₂₅, -(CH₂)₀₋₄-CO-(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-CO-(C₂-C₁₂ alkenyl), -(CH₂)₀₋₄-CO-(C₂-C₁₂ alkynyl), -(CH₂)₀₋₄-CO-(C₃-C₇ cycloalkyl), -(CH₂)₀₋₄-CO-aryl, -(CH₂)₀₋₄-CO-heteroaryl, -(CH₂)₀₋₄-CO-heterocyclyl, -(CH₂)₀₋₄-CO-O-R₂₁₅, -(CH₂)₀₋₄-SO₂-NR₂₂₀R₂₂₅, -(CH₂)₀₋₄-SO-(C₁-C₈ alkyl), -(CH₂)₀₋₄-SO₂-(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-SO₂-(C₃-C₇ cycloalkyl), -(CH₂)₀₋₄-N(H or R₂₁₅)-CO-O-R₂₁₅, -(CH₂)₀₋₄-N(H or R₂₁₅)-CO-N(R₂₁₅)₂, -(CH₂)₀₋₄-N-CS-N(R₂₁₅)₂, -(CH₂)₀₋₄-N(-H or R₂₁₅)-CO-R₂₂₀, -(CH₂)₀₋₄-NR₂₂₀R₂₂₅, -(CH₂)₀₋₄-O-CO-(C₁-C₆ alkyl), -(CH₂)₀₋₄-O-P(O)-(OR₂₄₀)₂, -(CH₂)₀₋₄-O-CO-N(R₂₁₅)₂, -(CH₂)₀₋₄-O-CS-N(R₂₁₅)₂,

$-(\text{CH}_2)_{0-4}-\text{O}-$ (R₂₁₅), $-(\text{CH}_2)_{0-4}-\text{O}-$ (R₂₁₅) $-\text{COOH}$, $-(\text{CH}_2)_{0-4}-\text{S}-$ (R₂₁₅),
 $-(\text{CH}_2)_{0-4}-\text{O}-$ (C₁-C₆ alkyl optionally substituted with 1, 2, 3,
or 5 -F), C₃-C₇ cycloalkyl, $-(\text{CH}_2)_{0-4}-\text{N}(\text{H}$ or R₂₁₅) $-\text{SO}_2-\text{R}_{220}$,

$-(\text{CH}_2)_{0-4}-\text{C}_3-\text{C}_7$ cycloalkyl, or
C₁-C₁₀ alkyl optionally substituted with 1, 2, or 3 R₂₀₅
groups, or

C₂-C₁₀ alkenyl or C₂-C₁₀ alkynyl, each of which is optionally
substituted with 1 or 2 R₂₀₅ groups, wherein

the aryl and heteroaryl groups at each occurrence are
optionally substituted with 1, 2, or 3 groups that are
independently R₂₀₅, R₂₁₀, or

C₁-C₆ alkyl substituted with 1, 2, or 3 groups that are
independently R₂₀₅ or R₂₁₀, and wherein

the heterocyclyl group at each occurrence is optionally
substituted with 1, 2, or 3 groups that are
independently R₂₁₀;

R₂₀₅ at each occurrence is independently selected from C₁-C₆
alkyl, halogen, -OH, -O-phenyl, -SH, -C≡N, -CF₃, C₁-C₆
alkoxy, NH₂, NH(C₁-C₆ alkyl) or N-(C₁-C₆ alkyl)(C₁-C₆ alkyl);

R₂₁₀ at each occurrence is independently selected from halogen,
C₁-C₆ alkoxy, C₁-C₆ haloalkoxy, -NR₂₂₀R₂₂₅, OH, C≡N, -CO-(C₁-C₄
alkyl), -SO₂-NR₂₃₅R₂₄₀, -CO-NR₂₃₅R₂₄₀, -SO₂-(C₁-C₄ alkyl), =O, or
C₁-C₆ alkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl or C₃-C₇ cycloalkyl,
each of which is optionally substituted with 1, 2, or 3
R₂₀₅ groups;

R₂₁₅ at each occurrence is independently selected from C₁-C₆
alkyl, $-(\text{CH}_2)_{0-2}-$ (aryl), C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₃-C₇
cycloalkyl, and $-(\text{CH}_2)_{0-2}-$ (heteroaryl), $-(\text{CH}_2)_{0-2}-$
(heterocyclyl), wherein

the aryl group at each occurrence is optionally substituted
with 1, 2, or 3 groups that are independently R₂₀₅ or
R₂₁₀, and wherein

the heterocyclyl and heteroaryl groups at each occurrence are optionally substituted with 1, 2, or 3 R₂₁₀;

R₂₂₀ and R₂₂₅ at each occurrence are independently selected from -H, -C₃-C₇ cycloalkyl, -(C₁-C₂ alkyl)-(C₃-C₇ cycloalkyl), -(C₁-C₆ alkyl)-O-(C₁-C₃ alkyl), -C₂-C₆ alkenyl, -C₂-C₆ alkynyl, -C₁-C₆ alkyl chain with one double bond and one triple bond, -aryl, -heteroaryl, and -heterocyclyl, or -C₁-C₁₀ alkyl optionally substituted with -OH, -NH₂ or halogen, wherein

the aryl, heterocyclyl and heteroaryl groups at each occurrence are optionally substituted with 1, 2, or 3 R₂₇₀ groups

R₂₃₅ and R₂₄₀ at each occurrence are independently H, or C₁-C₆ alkyl;

R₂₄₅ and R₂₅₀ at each occurrence are independently selected from -H, C₁-C₄ alkyl, C₁-C₄ alkylaryl, C₁-C₄ alkylheteroaryl, C₁-C₄ hydroxyalkyl, C₁-C₄ alkoxy, C₁-C₄ haloalkoxy, -(CH₂)₀₋₄-C₃-C₇ cycloalkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl, and phenyl; or

R₂₄₅ and R₂₅₀ are taken together with the carbon to which they are attached to form a carbocycle of 3, 4, 5, 6, or 7 carbon atoms, where one carbon atom is optionally replaced by a heteroatom selected from -O-, -S-, -SO₂-, and -NR₂₂₀-;

R₂₅₅ and R₂₆₀ at each occurrence are independently selected from -H, -(CH₂)₁₋₂-S(O)₀₋₂-(C₁-C₆ alkyl), -(C₁-C₄ alkyl)-aryl, -(C₁-C₄ alkyl)-heteroaryl, -(C₁-C₄ alkyl)-heterocyclyl, -aryl, -heteroaryl, -heterocyclyl, -(CH₂)₁₋₄-R₂₆₅-(CH₂)₀₋₄-aryl, -(CH₂)₁₋₄-R₂₆₅-(CH₂)₀₋₄-heteroaryl, -(CH₂)₁₋₄-R₂₆₅-(CH₂)₀₋₄-heterocyclyl, or

C₁-C₆ alkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl or -(CH₂)₀₋₄-C₃-C₇ cycloalkyl, each of which is optionally substituted with 1, 2, or 3 R₂₀₅ groups, wherein

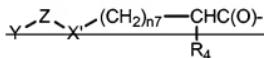
each aryl or phenyl is optionally substituted with 1, 2, or 3 groups that are independently R₂₀₅, R₂₁₀, or

C₁-C₆ alkyl substituted with 1, 2, or 3 groups that are independently R₂₀₅ or R₂₁₀, and wherein each heterocyclyl is optionally substituted with 1, 2, 3, or 4 R₂₁₀;

R₂₆₅ at each occurrence is independently -O-, -S- or -N(C₁-C₆ alkyl)-;

R₂₇₀ at each occurrence is independently R₂₀₅, halogen C₁-C₆ alkoxy, C₁-C₆ haloalkoxy, NR₂₃₅R₂₄₀, -OH, -C≡N, -CO-(C₁-C₄ alkyl), -SO₂-NR₂₃₅R₂₄₀, -CO-NR₂₃₅R₂₄₀, -SO₂-(C₁-C₄ alkyl), =O, or C₁-C₆ alkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl or -(CH₂)₀₋₄-C₃-C₇ cycloalkyl, each of which is optionally substituted with 1, 2, or 3 R₂₀₅ groups;

R_N is R'₁₀₀, -C(=O)-NR₁₀₀-R'₁₀₀, -C(=O)O-R'₁₀₀, -(CRR')₁₋₆R'₁₀₀, -C(=O)-(CRR')₀₋₆R₁₀₀, -C(=O)-(CRR')₁₋₆-O-R'₁₀₀, -C(=O)-(CRR')₁₋₆-S-R'₁₀₀, -C(=O)-(CRR')₁₋₆-C(=O)-R₁₀₀, -C(=O)-(CRR')₁₋₆-SO₂-R₁₀₀, or -C(=O)-(CRR')₁₋₆-NR₁₀₀-R'₁₀₀; ~~or~~



wherein

R₄ is selected from the group consisting of H, NH, NH-(CH₂)_{n6}-R₄, ~~+7-NHR₄, -NR₅₅C(O)R₄, C₁-C₄ alkyl-NHC(O)R₄, -(CH₂)_{n6}R₄, -O-C₁-C₄ alkenoyl, OH, C₁-C₆ aryloxy optionally substituted with 1, 2, or 3 groups that are independently halogen, C₁-C₄ alkyl, CO₂H, -C(O)-C₁-C₄ alkoxy, or C₁-C₄ alkoxy, C₁-C₆ alkoxy, aryl C₁-C₄ alkoxy, NR₅₅CO₂R₄₄, C₁-C₄ alkyl-NR₅₅CO₂R₄₄, -C≡N, -CF₃, -CF₂-CF₃, -C≡CH, -CH₂-CH=CH₂, -(CH₂)₁₋₄-R₄₋₁₇, -(CH₂)₄₋₁₇-NH-R₄₋₁₇, -O-(CH₂)_{n6}-R₄₋₁₇, -S-(CH₂)_{n6}-R₄₋₁₇, -(CH₂)₀₋₆-R₅₂, -(CH₂)₀₋₄-R₅₃-(CH₂)₀₋₄-R₅₄,~~

wherein

n₆ is 0, 1, 2, or 3,

n₇ is 0, 1, 2, or 3;

R_{4-1} is selected from the group consisting of $\text{SO}_2-(\text{C}_1-\text{C}_6$ alkyl), $\text{SO}(\text{C}_1-\text{C}_6$ alkyl), $\text{S}(\text{C}_1-\text{C}_6$ alkyl), $\text{S-CO}(\text{C}_1-\text{C}_6$ alkyl), $\text{SO}_2-\text{NR}_{4-2}\text{R}_{4-3}$, $\text{CO-C}_1-\text{C}_6$ alkyl, $\text{CO-NR}_{4-3}\text{R}_{4-4}$, and R_{4-2} and R_{4-3} are independently H, C_1-C_6 alkyl, or C_1-C_6 cycloalkyl;

R_{4-4} is alkyl, arylalkyl, alkanoyl, or arylalkanoyl;

R_{4-5} is H or C_1-C_6 alkyl;

R_5 is selected from the group consisting of C_1-C_6 cycloalkyl, C_1-C_6 alkyl optionally substituted with 1, 2, or 3 groups that are independently halogen, NR_6R_7 , C_1-C_6 alkoxy, C_1-C_6 heterocycloalkyl, C_1-C_6 heteroaryl, C_1-C_6 aryl, C_1-C_6 cycloalkyl C_1-C_6 alkyl, $\text{S-C}_1-\text{C}_6$ alkyl, $\text{SO}_2-\text{C}_1-\text{C}_6$ alkyl, $\text{-CO}_2\text{H}$, $\text{-CONR}_8\text{R}_9$, $\text{CO}_2-\text{C}_1-\text{C}_6$ alkyl, C_1-C_6 aryloxy, heteroaryl optionally substituted with 1, 2, or 3 groups that are independently C_1-C_6 alkyl, C_1-C_6 alkoxy, halogen, C_1-C_6 haloalkyl, or OH, heterocycloalkyl optionally substituted with 1, 2, or 3 groups that are independently C_1-C_6 alkyl, C_1-C_6 alkoxy, halogen, or C_1-C_6 alkanoyl, aryl optionally substituted with 1, 2, 3, or 4 groups that are independently halogen, OH, C_1-C_6 alkyl, C_1-C_6 alkoxy, or C_1-C_6 haloalkyl, and NR_6R_7 , wherein

R_6 and R_7 are independently selected from the group consisting of H, C_1-C_6 alkyl, C_1-C_6 alkanoyl, phenyl, $\text{SO}_2-\text{C}_1-\text{C}_6$ alkyl, phenyl C_1-C_6 alkyl,

R_8 is selected from the group consisting of SO_2 heteroaryl, -SO_2 -aryl, -SO_2 -heterocycloalkyl, $\text{-SO}_2-\text{C}_1-\text{C}_6$ alkyl, -C(O)NHR_9 , heterocycloalkyl, $\text{-S-C}_1-\text{C}_6$ alkyl, $\text{-S-C}_1-\text{C}_6$ alkanoyl, wherein

R_9 is aryl C_1-C_6 alkyl, C_1-C_6 alkyl, or H;

R_{50} is H or C_1-C_6 alkyl;

R_{51} is selected from the group consisting of aryl C_1-C_6 alkyl, C_1-C_6 alkyl optionally substituted with 1, 2, or

~~3 groups that are independently halogen, cyano, heteroaryl, NR₂R, C(O)NR₂R, C₁-C₄ cycloalkyl, or C₁-C₄ alkoxy; heterocycloalkyl optionally substituted with 1 or 2 groups that are independently C₁-C₄ alkyl, C₁-C₄ alkoxy, halogen, C₁-C₄ alkanoyl, aryl C₁-C₄ alkyl, and SO₂C₁-C₄ alkyl, alkenyl, alkynyl; heteroaryl optionally substituted with 1, 2, or 3 groups that are independently OH, C₁-C₄ alkyl, C₁-C₄ alkoxy, halogen, NH₂, NH(C₁-C₄ alkyl) or N(C₁-C₄ alkyl)(C₁-C₄ alkyl); heteroaryalkyl optionally substituted with 1, 2, or 3 groups that are independently C₁-C₄ alkyl, C₁-C₄ alkoxy, halogen, NH₂, NH(C₁-C₄ alkyl) or N(C₁-C₄ alkyl)(C₁-C₄ alkyl), aryl, heterocycloalkyl, C₁-C₄ cycloalkyl, and cycloalkylalkyl, wherein the aryl, heterocycloalkyl, C₁-C₄ cycloalkyl, and cycloalkylalkyl groups are optionally substituted with 1, 2, 3, 4 or 5 groups that are independently halogen, CN, NO₂, C₁-C₄ alkyl, C₁-C₄ alkoxy, C₁-C₄ alkanoyl, C₁-C₄ haloalkyl, C₁-C₄ haloalkoxy, hydroxy, C₁-C₄ hydroxyalkyl, C₁-C₄ alkoxy C₁-C₄ alkyl, C₁-C₄ thioalkoxy, C₁-C₄ thioalkoxy C₁-C₄ alkyl, or C₁-C₄ alkoxy C₁-C₄ alkoxy;~~

~~R₂₂ is heterocycloalkyl, heteroaryl, aryl, cycloalkyl, S(O)₂-C₁-C₄ alkyl, CO₂H, C(O)NH₂, C(O)NH(alkyl), C(O)N(alkyl)(alkyl), CO₂ alkyl, NHS(O)₂-C₁-C₄ alkyl, N(alkyl)S(O)₂-C₁-C₄ alkyl, S(O)₂-heteroaryl, S(O)₂-aryl, NH(arylalkyl), N(alkyl)(arylalkyl), thioalkoxy, or alkoxy, each of which is optionally substituted with 1, 2, 3, 4, or 5 groups that are independently alkyl, alkoxy, thioalkoxy, halogen, haloalkyl, haloalkoxy, alkanoyl, NO₂, CN, alkoxy carbonyl, or amino carbonyl,~~

R_{53} is absent, —O—, —C(O)—, —NH—, —N(alkyl)–, —NH—S(O)₀₋₂—, —N(alkyl)–S(O)₀₋₂—, —S(O)₀₋₂—NH—, —S(O)₀₋₂—N(alkyl)–, —NH—C(S)–, or —N(alkyl)–C(S)–,

R_{54} is heteroaryl, aryl, arylalkyl, heterocycloalkyl, CO₂H, —CO₂–alkyl, —C(O)NH(alkyl), —C(O)N(alkyl)–(alkyl), —C(O)NH₂, C₁–C₄–alkyl, OH, aryloxy, alkoxy, arylalkoxy, NH₂, NH(alkyl), N(alkyl)–(alkyl), or C₁–C₄–alkyl—CO₂–C₁–C₄–alkyl, each of which is optionally substituted with 1, 2, 3, 4, or 5 groups that are independently alkyl, alkoxy, CO₂H, CO₂–alkyl, thioalkoxy, halogen, haloalkyl, haloalkoxy, hydroxylalkyl, alkanyl, NO₂, CN, alkoxy carbonyl, or aminocarbonyl,

X' is selected from the group consisting of C₁–C₄–alkylideneyl optionally substituted with 1, 2, or 3 methyl groups; and NR₄₋₆, or

R_4 and R_{4-6} combine to form (CH₂)_{n+1}, wherein n is 1, 2, 3, or 4;

Z is selected from the group consisting of a bond, SO₂, SO, S⁺ and C(O)–,

Y is selected from the group consisting of H, C₁–C₄–haloalkyl, C₁–C₄–heterocycloalkyl, C₁–C₄–aryl, OH, —N(Y₁)(Y₂), C₁–C₁₀–alkyl optionally substituted with 1 thru 3 substituents which can be the same or different and are selected from the group consisting of halogen, hydroxy, alkoxy, thioalkoxy, and haloalkoxy, C₁–C₄–cycloalkyl optionally substituted with 1, 2, or 3 groups independently selected from C₁–C₄–alkyl, and halogen; alkoxy, aryl optionally substituted with halogen, alkyl, alkoxy, CN or NO₂; arylalkyl optionally substituted with halogen, alkyl, alkoxy, CN or NO₂; wherein

Y₁ and Y₂ are the same or different and are H, C₁–C₄–alkyl optionally substituted with 1, 2, or 3 substituents selected from the group consisting of halogen, C₁–C₄–alkoxy, C₁–C₄–cycloalkyl, and OH; C₂–C₆–alkenyl; C₂–C₆–

~~alkanoyl, phenyl-, $\text{SO}_2\text{-C}_6\text{-C}_4$ -alkyl, phenyl- $\text{C}_6\text{-C}_4$ -alkyl,
or $\text{C}_6\text{-C}_6$ -cycloalkyl- $\text{C}_6\text{-C}_4$ -alkyl, or
 Y_1 , Y_2 , and the nitrogen to which they are attached form a
ring selected from the group consisting of piperazinyl,
piperidinyl, morpholinyl, and pyrrolidinyl, wherein each
ring is optionally substituted with 1, 2, 3, or 4
groups that are independently $\text{C}_6\text{-C}_6$ -alkyl, $\text{C}_6\text{-C}_6$ -alkoxy,
 $\text{C}_6\text{-C}_6$ -alkoxy- $\text{C}_6\text{-C}_6$ -alkyl, or halogen,~~

R₁₀₀ and R'₁₀₀ independently represent aryl, heteroaryl, -aryl-W-
aryl, -aryl-W-heteroaryl, -aryl-W-heterocyclyl, -heteroaryl-
W-aryl, -heteroaryl-W-heteroaryl, -heteroaryl-W-
heterocyclyl, -heterocyclyl-W-aryl, -heterocyclyl-W-
heteroaryl, -heterocyclyl-W-heterocyclyl, - $\text{CH}[(\text{CH}_2)_{0-2}\text{-O-}$
R₁₅₀]- $(\text{CH}_2)_{0-2}$ -aryl, - $\text{CH}[(\text{CH}_2)_{0-2}\text{-O-}$ R₁₅₀]- $(\text{CH}_2)_{0-2}$ -heterocyclyl or
- $\text{CH}[(\text{CH}_2)_{0-2}\text{-O-}$ R₁₅₀]- $(\text{CH}_2)_{0-2}$ -heteroaryl, where the ring
portions of each are optionally substituted with 1, 2, or 3
groups independently selected from

-OR, -NO₂, halogen, -C≡N, -OCF₃, -CF₃, -(CH₂)₀₋₄O-
P(=O)(OR')(OR''), -(CH₂)₀₋₄-CO-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄O-(CH₂)₀₋₄-
CONR₁₀₂R₁₀₂', -(CH₂)₀₋₄-CO-(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-CO-(C₂-
C₁₂ alkenyl), -(CH₂)₀₋₄-CO-(C₂-C₁₂ alkynyl), -(CH₂)₀₋₄-CO-
(CH₂)₀₋₄(C₃-C₇ cycloalkyl), -(CH₂)₀₋₄-R₁₁₀, -(CH₂)₀₋₄-R₁₂₀,
-(CH₂)₀₋₄-R₁₃₀, -(CH₂)₀₋₄-CO-R₁₁₀, -(CH₂)₀₋₄-CO-R₁₂₀, -(CH₂)₀₋₄-
CO-R₁₃₀, -(CH₂)₀₋₄-CO-R₁₄₀, -(CH₂)₀₋₄-CO-O-R₁₅₀, -(CH₂)₀₋₄-
SO₂-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄-SO-(C₁-C₈ alkyl), -(CH₂)₀₋₄-SO₂-
(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-SO₂-(CH₂)₀₋₄-(C₃-C₇ cycloalkyl),
-(CH₂)₀₋₄-N(R₁₅₀)-CO-O-R₁₅₀, -(CH₂)₀₋₄-N(R₁₅₀)-CO-N(R₁₅₀)₂,
-(CH₂)₀₋₄-N(R₁₅₀)-CS-N(R₁₅₀)₂, -(CH₂)₀₋₄-N(R₁₅₀)-CO-R₁₀₅,
-(CH₂)₀₋₄-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄-R₁₄₀, -(CH₂)₀₋₄-O-CO-(C₁-C₆
alkyl), -(CH₂)₀₋₄-O-P(O)-(O-R₁₁₀)₂, -(CH₂)₀₋₄-O-CO-N(R₁₅₀)₂,
-(CH₂)₀₋₄-O-CS-N(R₁₅₀)₂, -(CH₂)₀₋₄-O-(R₁₅₀), -(CH₂)₀₋₄-O-
R₁₅₀'-COOH, -(CH₂)₀₋₄S-(R₁₅₀), -(CH₂)₀₋₄-N(R₁₅₀)-SO₂-R₁₀₅,

- (CH₂)₀₋₄₋ C₃-C₇ cycloalkyl, (C₂-C₁₀) alkenyl, or (C₂-C₁₀) alkynyl, or

R₁₀₀ is C₁-C₁₀ alkyl optionally substituted with 1, 2, or 3 R₁₁₅ groups, or

R₁₀₀ is -(C₁-C₆ alkyl)-O-C₁-C₆ alkyl) or -(C₁-C₆ alkyl)-S-(C₁-C₆ alkyl), each of which is optionally substituted with 1, 2, or 3 R₁₁₅ groups, or

R₁₀₀ is C₃-C₈ cycloalkyl optionally substituted with 1, 2, or 3 R₁₁₅ groups;

W is -(CH₂)₀₋₄₋, -O-, -S(O)₀₋₂₋, -N(R₁₃₅)-, -CR(OH)- or -C(O)-;

R₁₀₂ and R_{102'} independently are hydrogen, or

C₁-C₁₀ alkyl optionally substituted with 1, 2, or 3 groups that are independently halogen, aryl or -R₁₁₀;

R₁₀₅ and R'₁₀₅ independently represent -H, -R₁₁₀, -R₁₂₀, C₃-C₇ cycloalkyl, -(C₁-C₂ alkyl)-(C₃-C₇ cycloalkyl), -(C₁-C₆ alkyl)-O-(C₁-C₃ alkyl), C₂-C₆ alkenyl, C₂-C₆ alkynyl, or C₁-C₆ alkyl chain with one double bond and one triple bond, or

C₁-C₆ alkyl optionally substituted with -OH or -NH₂; or,

C₁-C₆ alkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, or

R₁₀₅ and R'₁₀₅ together with the atom to which they are attached form a 3 to 7 membered carbocyclic ring, where one member is optionally a heteratom selected from -O-, -S(O)₀₋₂₋, -N(R₁₃₅)-, the ring being optionally substituted with 1, 2 or three R₁₄₀ groups;

R₁₁₅ at each occurrence is independently halogen, -OH, -CO₂R₁₀₂, -C₁-C₆ thioalkoxy, -CO₂-phenyl, -NR₁₀₅R'₁₃₅, -SO₂-(C₁-C₈ alkyl), -C(=O)R₁₈₀, R₁₈₀, -CONR₁₀₅R'₁₀₅, -SO₂NR₁₀₅R'₁₀₅, -NH-CO-(C₁-C₆ alkyl), -NH-C(=O)-OH, -NH-C(=O)-OR, -NH-C(=O)-O-phenyl, -O-C(=O)-(C₁-C₆ alkyl), -O-C(=O)-amino, -O-C(=O)-mono- or dialkylamino, -O-C(=O)-phenyl, -O-(C₁-C₆ alkyl)-CO₂H, -NH-SO₂-(C₁-C₆ alkyl), C₁-C₆ alkoxy or C₁-C₆ haloalkoxy;

R₁₃₅ is C₁-C₆ alkyl, C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₃-C₇ cycloalkyl, -(CH₂)₀₋₂₋(aryl), -(CH₂)₀₋₂₋(heteroaryl), or -(CH₂)₀₋₂₋(heterocyclyl);

R₁₄₀ is heterocyclyl optionally substituted with 1, 2, 3, or 4 groups independently selected from C₁-C₆ alkyl, C₁-C₆ alkoxy, halogen, hydroxy, cyano, nitro, amino, mono(C₁-C₆)alkylamino, di(C₁-C₆)alkylamino, C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₁-C₆ haloalkyl, C₁-C₆ haloalkoxy, amino(C₁-C₆)alkyl, mono(C₁-C₆)alkylamino(C₁-C₆)alkyl, di(C₁-C₆)alkylamino(C₁-C₆)alkyl, and =O;

R₁₅₀ is hydrogen, C₃-C₇ cycloalkyl, -(C₁-C₂ alkyl)-(C₃-C₇ cycloalkyl), C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₁-C₆ alkyl with one double bond and one triple bond, -R₁₁₀, -R₁₂₀, or C₁-C₆ alkyl optionally substituted with 1, 2, 3, or 4 groups independently selected from -OH, -NH₂, C₁-C₃ alkoxy, R₁₁₀, and halogen;

R_{150'} is C₃-C₇ cycloalkyl, -(C₁-C₃ alkyl)-(C₃-C₇ cycloalkyl), C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₁-C₆ alkyl with one double bond and one triple bond, -R₁₁₀, -R₁₂₀, or C₁-C₆ alkyl optionally substituted with 1, 2, 3, or 4 groups independently selected from -OH, -NH₂, C₁-C₃ alkoxy, R₁₁₀, and halogen;

R₁₈₀ is selected from morpholinyl, thiomorpholinyl, piperazinyl, piperidinyl, homomorpholinyl, homothiomorpholinyl, homothiomorpholinyl S-oxide, homothiomorpholinyl S,S-dioxide, pyrrolinyl and pyrrolidinyl, each of which is optionally substituted with 1, 2, 3, or 4 groups independently selected from C₁-C₆ alkyl, C₁-C₆ alkoxy, halogen, hydroxy, cyano, nitro, amino, mono(C₁-C₆)alkylamino, di(C₁-C₆)alkylamino, C₂-C₆ alkenyl, C₂-C₆ alkynyl, C₁-C₆ haloalkyl, C₁-C₆ haloalkoxy, amino(C₁-C₆)alkyl, mono(C₁-C₆)alkylamino(C₁-C₆)alkyl, di(C₁-C₆)alkylamino(C₁-C₆)alkyl, and =O;

R_{110} is aryl optionally substituted with 1 or 2 R_{125} groups;
 R_{125} at each occurrence is independently halogen, amino, mono- or dialkylamino, -OH, -C≡N, -SO₂-NH₂, -SO₂-NH-C₁-C₆ alkyl, -SO₂-N(C₁-C₆ alkyl)₂, -SO₂-(C₁-C₄ alkyl), -CO-NH₂, -CO-NH-C₁-C₆ alkyl, or -CO-N(C₁-C₆ alkyl)₂, or C₁-C₆ alkyl, C₂-C₆ alkenyl or C₂-C₆ alkynyl, each of which is optionally substituted with 1, 2, or 3 groups that are independently selected from C₁-C₃ alkyl, halogen, -OH, -SH, -C≡N, -CF₃, C₁-C₃ alkoxy, amino, and mono- and dialkylamino, or C₁-C₆ alkoxy optionally substituted with one, two or three of halogen;

R_{120} is heteroaryl, which is optionally substituted with 1 or 2 R_{125} groups; and

R_{130} is heterocyclyl optionally substituted with 1 or 2 R_{125} groups; and

R_2 is selected from the group consisting of H; C₁-C₆ alkyl, optionally substituted with 1, 2, or 3 substituents that are independently selected from the group consisting of C₁-C₂ alkyl, halogen, OH, SH, C≡N, CF₃, C₁-C₃ alkoxy, and NR₁₋₂, wherein

R_{1-2} and R_{1-3} are H or C₁-C₆ alkyl,

(CH₂)₀₋₄-aryl, (CH₂)₀₋₄-heteroaryl, C₂-C₆ alkenyl, C₂-C₆ alkynyl, CONR₁₋₂R₁₋₃, SO₂NR₁₋₂R₁₋₃, CO₂H, and CO₂-(C₁-C₄ alkyl),

R₃ is selected from the group consisting of H, C₁-C₄ alkyl, optionally substituted with 1, 2, or 3 substituents independently selected from the group consisting of e-e. alkyl, halogen, OH, SH, CN, CF₃, C₁-C₄ alkoxy, and NR₁-R₂, -(CH₂)₀₋₄-aryl, -(CH₂)₀₋₄-heteroaryl, C₁-C₄ alkenyl, C₁-C₄ alkynyl, CO-NR₁-R₂, SO₂-NR₁-R₂, CO₂H, and CO-O-(C₁-C₄ alkyl);

wherein

R₁ and R₂ at each occurrence are independently selected from the group consisting of C₁-C₄ alkyl optionally substituted with 1, 2, or 3 groups independently selected from the group consisting of -OH, NH₂, phenyl and halogen, C₁-C₄ cycloalkyl, (C₁-C₄ alkyl) (C₁-C₄ cycloalkyl), (C₁-C₄ alkyl) O (C₁-C₄ alkyl), C₁-C₄ alkenyl, C₁-C₄ alkynyl, C₁-C₄ alkyl chain with one double bond and one triple bond, aryl, heteroaryl, heterocycloalkyl, or

R₁, R₂ and the nitrogen to which they are attached form a 5, 6, or 7 membered heterocycloalkyl or heteroaryl group, wherein said heterocycloalkyl or heteroaryl group is optionally fused to a benzene, pyridine, or pyrimidine ring, and said groups are unsubstituted or substituted with 1, 2, 3, 4, or 5 groups that at each occurrence are independently C₁-C₄ alkyl, C₁-C₄ alkoxy, halogen, halo C₁-C₄ alkyl, halo C₁-C₄ alkoxy, CN, NO₂, NH₂, NH(C₁-C₄ alkyl), N(C₁-C₄ alkyl) (C₁-C₄ alkyl), -OH, C(O)NH₂, C(O)NH(C₁-C₄ alkyl), C(O)N(C₁-C₄ alkyl) (C₁-C₄ alkyl), C₁-C₄ alkoxy C₁-C₄ alkyl, C₁-C₄ thioalkoxy, and C₁-C₄ thioalkoxy C₁-C₄ alkyl,

or wherein,

R₁, R₂ and the carbon to which they are attached form a carbocycle of three thru seven carbon atoms, wherein one carbon

atom is optionally replaced by a group selected from O , S ,
 -SO_2 , or NR_{1-2} .

2-4. (Canceled)

5. (Currently Amended) A compound according to claim 4_1, wherein R_1 is

$\text{-C}_1\text{-C}_6$ alkyl-aryl, $\text{-C}_1\text{-C}_6$ alkyl-heteroaryl, or $\text{-C}_1\text{-C}_6$ alkyl-heterocyclyl, where the ring portions of each are optionally substituted with 1, 2, 3, or 4 groups independently selected from halogen, -OH , -SH , $\text{-C}\equiv\text{N}$, -NO_2 , $\text{-NR}_{105}\text{R'}_{105}$, $\text{-CO}_2\text{R}$, $\text{-N}(\text{R})\text{COR}'$, or $\text{-N}(\text{R})\text{SO}_2\text{R}'$ (where R_{105} , R'_{105} , R and R' are as defined above), $\text{-C}(\text{=O})-(\text{C}_1\text{-C}_4)$ alkyl, -SO_2 -amino, -SO_2 -mono or dialkylamino, $\text{-C}(\text{=O})$ -amino, $\text{-C}(\text{=O})$ -mono or dialkylamino, $\text{-SO}_2-(\text{C}_1\text{-C}_4)$ alkyl, or $\text{C}_1\text{-C}_6$ alkoxy optionally substituted with 1, 2, or 3 groups which are independently selected from halogen, or

$\text{C}_1\text{-C}_6$ cycloalkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, -OH , -SH , $\text{-C}\equiv\text{N}$, -CF_3 , $\text{C}_1\text{-C}_6$ alkoxy, amino, $\text{C}_1\text{-C}_6$ alkyl and mono or dialkylamino, or

$\text{C}_1\text{-C}_6$ alkyl optionally substituted with 1, 2, or 3 groups independently selected from halogen, -OH , -SH , $\text{-C}\equiv\text{N}$, -CF_3 , $\text{C}_1\text{-C}_6$ alkoxy, amino, mono or dialkylamino and $\text{C}_1\text{-C}_6$ alkyl, or

$\text{C}_1\text{-C}_6$ alkenyl or $\text{C}_2\text{-C}_{10}$ alkynyl each of which is optionally substituted with 1, 2, or 3 groups independently selected from halogen, -OH , -SH , $\text{-C}\equiv\text{N}$, -CF_3 , $\text{C}_1\text{-C}_6$ alkoxy, amino, $\text{C}_1\text{-C}_6$ alkyl and mono or dialkylamino; and the heterocyclyl group is optionally further substituted with ~~exo~~.

6. (Original) A compound according to claim 1 wherein:
 R_N is $\text{-C}(\text{=O})-\text{R}_{100}$; and

R_{100} represents aryl, or heteroaryl, where the ring portions of each are optionally substituted with 1, 2, or 3 groups independently selected from

-OR, -NO₂, C₁-C₆ alkyl, halogen, -C≡N, -OCF₃, -CF₃, -(CH₂)₀₋₄-O-P(=O)(OR)(OR'), -(CH₂)₀₋₄-CO-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄-O-(CH₂)₀₋₄-CONR₁₀₂R'₁₀₂, -(CH₂)₀₋₄-CO-(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-CO-(C₂-C₁₂ alkenyl), -(CH₂)₀₋₄-CO-(C₂-C₁₂ alkynyl), -(CH₂)₀₋₄-CO-(CH₂)₀₋₄(C₃-C₇ cycloalkyl), -(CH₂)₀₋₄-R₁₁₀, -(CH₂)₀₋₄-R₁₂₀, -(CH₂)₀₋₄-R₁₃₀, -(CH₂)₀₋₄-CO-R₁₁₀, -(CH₂)₀₋₄-CO-R₁₂₀, -(CH₂)₀₋₄-CO-R₁₃₀, -(CH₂)₀₋₄-CO-R₁₄₀, -(CH₂)₀₋₄-CO-O-R₁₅₀, -(CH₂)₀₋₄-SO₂-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄-SO-(C₁-C₈ alkyl), -(CH₂)₀₋₄-SO₂-(C₁-C₁₂ alkyl), -(CH₂)₀₋₄-SO₂-(CH₂)₀₋₄-(C₃-C₇ cycloalkyl), -(CH₂)₀₋₄-N(R₁₅₀)-CO-O-R₁₅₀, -(CH₂)₀₋₄-N(R₁₅₀)-CO-N(R₁₅₀)₂, -(CH₂)₀₋₄-N(R₁₅₀)-CS-N(R₁₅₀)₂, -(CH₂)₀₋₄-N(R₁₅₀)-CO-R₁₀₅, -(CH₂)₀₋₄-NR₁₀₅R'₁₀₅, -(CH₂)₀₋₄-R₁₄₀, -(CH₂)₀₋₄-O-CO-(C₁-C₆ alkyl), -(CH₂)₀₋₄-O-P(O)-(O-R₁₁₀)₂, -(CH₂)₀₋₄-O-CO-N(R₁₅₀)₂, -(CH₂)₀₋₄-O-CS-N(R₁₅₀)₂, -(CH₂)₀₋₄-O-(R₁₅₀), -(CH₂)₀₋₄-O-R₁₅₀'-COOH, -(CH₂)₀₋₄-S-(R₁₅₀), -(CH₂)₀₋₄-N(R₁₅₀)-SO₂-R₁₀₅, -(CH₂)₀₋₄-(C₃-C₇ cycloalkyl, (C₂-C₁₀) alkenyl, or (C₂-C₁₀) alkynyl.

7. (Currently Amended) A compound according to claim 1 wherein:

R_C is hydrogen, -(CR₂₄₅R₂₅₀)₀₋₄-aryl, -(CR₂₄₅R₂₅₀)₀₋₄-heteroaryl, -(CR₂₄₅R₂₅₀)₀₋₄-heterocyclyl,
 ~~C_6-C_6 alkyl optionally substituted with 1, 2, or 3 groups independently selected from the group consisting of R₂₀₅, R₁₁₀, R₁₂₀, R₁₃₀, OC=ONR₂₃₅R₂₄₀, S(=O)₀₋₂(C₆-C₆ alkyl), -SH, and -S(=O)₂NR₂₃₅R₂₄₀,~~
~~-(CH₂)₀₋₂-(C₆-C₆) cycloalkyl wherein the cycloalkyl is optionally substituted with 1, 2, or 3 groups independently selected from the group consisting of R₂₀₅, CO₂H, and CO₂-(C₄-C₄ alkyl), or~~

~~C₂-C₄-alkenyl or C₂-C₄-alkynyl, each of which is optionally substituted with 1, 2, or 3 independently selected R₂₀₀ groups,~~ wherein

each aryl and heteroaryl is optionally substituted with 1, 2, or 3 R₂₀₀, and wherein each heterocyclyl is optionally substituted with 1, 2, 3, or 4 independently selected R₂₁₀ groups.

8. (Canceled)

9. (Previously Presented) A compound according to claim 1 selected from the group consisting of:

N-(3,5-difluorobenzyl)-N-{(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}-5-methyl-N',N'-dipropylisophthalamide;

N-[2-(3,5-difluorophenyl)ethyl]-N-{(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}-5-methyl-N',N'-dipropylisophthalamide;

3-[(2-(3,5-difluorophenyl)ethyl){(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}amino]sulfonyl]-N,N-dipropylbenzamide;

N-(3,5-difluorobenzyl)-N-((2R)-3-{{(4R)-6-ethyl-2,2-dioxido-3,4-dihydro-1H-isothiochromen-4-yl]amino}-2-hydroxypropyl)-5-methyl-N',N'-dipropylisophthalamide;

N-[2-(3,5-difluorophenyl)ethyl]-N-((2R)-3-{{(4R)-6-ethyl-2,2-dioxido-3,4-dihydro-1H-isothiochromen-4-yl]amino}-2-hydroxypropyl)-5-methyl-N',N'-dipropylisophthalamide;

3-{{[2-(3,5-difluorophenyl)ethyl]{(2R)-3-{{(4R)-6-ethyl-2,2-dioxido-3,4-dihydro-1H-isothiochromen-4-yl]amino}-2-hydroxypropyl}amino}sulfonyl}-N,N-dipropylbenzamide;

N-(3,5-difluorobenzyl)-N-{(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}-N',N',5-trimethylisophthalamide;

N-[2-(3,5-difluorophenyl)ethyl]-N-{(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}-N',N',5-trimethylisophthalamide;

3-{{[2-(3,5-difluorophenyl)ethyl]{(2R)-2-hydroxy-3-[(3-iodobenzyl)amino]propyl}amino}sulfonyl}-N,N-dimethylbenzamide;

N-(3,5-difluorobenzyl)-*N*-((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)-*N'*,*N'*,5-trimethylisophthalamide;

N-[2-(3,5-difluorophenyl)ethyl]-*N*-((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)-*N'*,*N'*,5-trimethylisophthalamide;

3-{{[2-(3,5-difluorophenyl)ethyl]}((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)amino}sulfonyl)-*N*,*N*-dimethylbenzamide;

N-(3-chloro-5-fluorobenzyl)-*N*-{(2*R*)-2-hydroxy-3-[{(3-iodobenzyl)amino}propyl]-5-methyl-*N'*,*N'*-dipropylisophthalamide;

N-[2-(3-chloro-5-fluorophenyl)ethyl]-*N*-{(2*R*)-2-hydroxy-3-[{(3-iodobenzyl)amino}propyl]-5-methyl-*N'*,*N'*-dipropylisophthalamide;

3-{{[2-(3-chloro-5-fluorophenyl)ethyl]}((2*R*)-2-hydroxy-3-[{(3-iodobenzyl)amino}propyl]amino)sulfonyl)-*N*,*N*-dipropylbenzamide;

N-(3-chloro-5-fluorobenzyl)-*N*-((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)-5-methyl-*N'*,*N'*-dipropylisophthalamide;

N-[2-(3-chloro-5-fluorophenyl)ethyl]-*N*-((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)-5-methyl-*N'*,*N'*-dipropylisophthalamide;

3-{{[2-(3-chloro-5-fluorophenyl)ethyl]}((2*R*)-3-{[(4*R*)-6-ethyl-2,2-dioxido-3,4-dihydro-1*H*-isothiochromen-4-yl]amino}-2-hydroxypropyl)amino}sulfonyl)-*N*,*N*-dipropylbenzamide;

N-[(2*R*)-3-(benzylamino)-2-hydroxypropyl]-*N*-(3,5-difluorobenzyl)-5-methyl-*N'*,*N'*-dipropylisophthalamide;

N-[(2*R*)-3-(benzylamino)-2-hydroxypropyl]-*N*-[2-(3,5-difluorophenyl)ethyl]-5-methyl-*N'*,*N'*-dipropylisophthalamide;

3-{{[(2*R*)-3-(benzylamino)-2-hydroxypropyl]}[2-(3,5-difluorophenyl)ethyl]amino}sulfonyl)-*N*,*N*-dipropylbenzamide; and salts thereof.

10. (Previously Presented) A pharmaceutical composition comprising a compound according to claim 1, in combination with a physiologically acceptable carrier or excipient.

11-12. (Canceled)

13. (Withdrawn) A method for treating a patient who has, or in preventing a patient from getting, a disease or condition selected from the group consisting of Alzheimer's disease, for helping prevent or delay the onset of Alzheimer's disease, for treating patients with mild cognitive impairment (MCI) and preventing or delaying the onset of Alzheimer's disease in those who would progress from MCI to AD, for treating Down's syndrome, for treating humans who have Hereditary Cerebral Hemorrhage with Amyloidosis of the Dutch-Type, for treating cerebral amyloid angiopathy and preventing its potential consequences, i.e. single and recurrent lobar hemorrhages, for treating other degenerative dementias, including dementias of mixed vascular and degenerative origin, dementia associated with Parkinson's disease, dementia associated with progressive supranuclear palsy, dementia associated with cortical basal degeneration, or diffuse Lewy body type of Alzheimer's disease and who is in need of such treatment, comprising administering to such patient a therapeutically effective amount of a compound of claim 1.

14. (Withdrawn) A method for the treatment or prevention of Alzheimer's disease, mild cognitive impairment Down's syndrome, Hereditary Cerebral Hemorrhage with Amyloidosis of the Dutch-Type, cerebral amyloid angiopathy, other degenerative dementias, dementias of mixed vascular and degenerative origin, dementia associated with Parkinson's disease, dementia associated with progressive supranuclear palsy, dementia associated with cortical

basal degeneration, diffuse Lewy body type of Alzheimer's disease comprising administration of a therapeutically effective amount of a compound or salt according to Claim 1, to a patient in need thereof.

15. (Withdrawn) A method for making a compound of claim 1.